

EXPERIMENTAL DESIGN FOR COMPLEX SYSTEMS

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5 BACKGROUND OF THE INVENTION**1. Field of the Invention**

This invention relates generally to the manufacture of high performance semiconductor devices. More specifically, this invention relates to a systematic approach to an experimental design for large, complex systems. Even more specifically, this
10 invention relates to a method for systematically designing experiments when prior knowledge of the many factor and multiple responses is expressed as a network of cause-effect relationships.

2. Discussion of the Related Art

The design of a new semiconductor device and the process for manufacturing the
15 new semiconductor device has three phases: the development phase during which processing alternatives are still under evaluation and the nominal process targets continue to be tuned; the pre-production phase during which the process targets are more-or-less set, processing experience is acquired, and appropriate tolerance windows are determined; and the production phase during which both the process target and tolerance windows are
20 more-or-less fixed, and the full resources of the manufacturing line are committed in volume. The present invention focuses on the pre-production phase during which process targets are substantially set, but process experience is to be accumulated, and during which appropriate tolerance windows need to be determined.

The pre-production phase plays an essential role in managing the manufacturing
25 risk factor, providing a time period for determining problematic and challenging process steps, for investigating the range over which product can be manufactured successfully, and for reliability stress testing. The scale of current semiconductor manufacturing processes magnifies all of the issues, for example, contemporary semiconductor processes have 300-400 value added steps, any of which is a source of poor quality and/or reliability.
30 A cost of a single test batch can exceed one quarter million dollars. A delay in market

entry equal to one cycle of learning (the production time required to make one batch) approaches two orders of magnitude more.

Competitive pressures have provided strong incentives to keep pre-production costs to an effective minimum. In the statistics literature, it is well recognized that

5 appropriately designed experiments more fully characterize processes than, for example, multiple repetitions of the nominal process. The pre-production assessment of a new semiconductor manufacturing process typically involves many factors and can be from 30 – 50 and sometimes more, multiple blocks, typically from 5 – 15 blocks, and several responses, typically from 3 – 8 responses. One practical constraint is for each

10 experimental block to be self-contained, in the sense that each block supports an analysis without necessarily requiring results from other blocks. A complementary goal has the entire ensemble of experimental blocks covering the process space well. Subject matter expertise is both available and desirable and can be organized as a network of likely cause-effect relationships.

15 The present invention thus presents a systematic approach to the pre-production problem, including objectives, constraints, overarching model, blocking structures, split and skew factors and self-containing blocks.

SUMMARY OF THE INVENTION

According to the present invention, the foregoing and other objects and

20 advantages are achieved a systematic approach to forming experimental designs for large, complex systems after an idea for a product is formed. In accordance with a first aspect of the invention, critical variables for the product are determined by experts in the field, a design matrix U_k is defined, a base design matrix X is generated, $Y(P) = (I - B(B^T B)^{-1} B^T)(X P) / U$ & Wynn's criterion is defined, where P is a permutation matrix, I is an identity

25 matrix, B is a blocking matrix, B^T is a transposed matrix of B and A is a matrix composed of causal map-base coefficients and wherein a design matrix U_k is created. The index $k \leftarrow k+1$ is set and an algorithm to choose the best of random column permutation matrices P and an algorithm to choose the best column permutation matrix P that is near a previous solution and setting $U_k \leftarrow [X P^k \text{ with rows from } U_{k-1} \text{ appended}]$.

30 In accordance with another aspect of the invention, it determined whether the design U_k is large enough and if not the process described above is repeated until the

design U_k is large enough. If it is determined that the design U_k is large enough prototype products are manufactured, model responses are determined from the prototype wafers and determining whether the model responses are adequate.

5 In another aspect of the invention, if the model responses are adequate, tolerances for the product are assessed and proposed. If the tolerances assessed and proposed are manufacturable, the product is passed to full-scale production.

In another aspect of the invention, if the model responses are not adequate, the experimental design is repeated to create further U_k . This procedure is repeated until a design U_k is achieved that indicates that the model is adequate.

10 In another aspect of the invention, if the tolerances assessed and proposed are not manufacturable, the design experiment is repeated until a design U_k is achieved that provides a manufacturable product.

15 The described invention thus provides a method for a systematic approach to forming experimental designs for large, complex systems after an idea for a product is formed.

20 The present invention is better understood upon consideration of the detailed description below, in conjunction with the accompanying drawings. As will become readily apparent to those skilled in the art from the following description, there is shown and described an embodiment of this invention simply by way of illustration of the best mode to carry out the invention. As will be realized, the invention is capable of other embodiments and its several details are capable of modifications in various obvious aspects, all without departing from the scope of the invention. Accordingly, the drawings and detailed description will be regarded as illustrative in nature and not as restrictive.

BRIEF DESCRIPTION OF THE DRAWINGS

25 The novel features believed characteristic of the invention are set forth in the appended claims. The invention itself, however, as well as a preferred mode of use, and further objects and advantages thereof, will best be understood by reference to the following detailed description of an illustrative embodiment when read in conjunction with the accompanying drawings, wherein:

Figure 1 is a flow diagram showing the steps for the development of a semiconductor device from the “Idea for a Product” to the “Production” of the semiconductor product;

Figure 2 is flow diagram showing the steps for the pre-production phase of the development of a semiconductor product;

Figure 3 is a diagram of a Causal-Effect Diagram;

Figures 3A – 3G are sub-diagrams associated with the Causal – Effect Diagram shown in **Figure 3**; and

Figure 4 is a causal map derived from information contained in the Causal-Effect Diagram shown in **Figure 3**.

DETAILED DESCRIPTION

Reference is now made in detail to a specific embodiment or specific embodiments of the present invention that illustrate the best mode or modes presently contemplated by the inventors for practicing the invention.

Figure 1 is a flow diagram **100** showing the steps for the production of a semiconductor device from the initial “idea” for the semiconductor device through the production of the semiconductor device. The first step is the generation of an “Idea for a Product” **Figure 1, Step A, 102**. The next step is the “Initial/Further Development of Idea” that comprehends gathering data associated with the desired product, **Figure 1, Step B, 104**. The next step is to “Design/Refine Experiment” **Figure 1, Step C, 106** and as indicated by the arrow **108** the flow goes to **Figure 2 Step A**. **Figure 2** is a flow diagram **200** showing the steps for the pre-production phase of the present invention and will be discussed below. After the pre-production phase shown in **Figure 2** is completed the flow returns to the flow diagram **100** as shown by arrow **110**. Prototype wafers manufactured using designed experimental values are evaluated at **112**. The responses from the prototype wafers are empirically measured and mathematically modeled at **114**. It is determined at **116** whether the response models are adequate. If the response models are not adequate, the flow returns to **Figure 2, Step I** as indicated by arrow **118**. If the models are adequate, the next step at **120** is to assess and propose tolerances. It is determined at **122** if the proposed tolerances can be manufactured. If the proposed tolerances can be

manufactured, the designed wafers are sent into production at **124**. If the proposed tolerances cannot be manufactured, the flow is returned to **Step B** as indicated by the arrow **126**.

Figure 2 is a flow diagram **200** showing the steps of the pre-production phase of the present invention. Arrow **202** indicates the flow from **Step C, Figure 1**. Referring to **Figure 2, Step A**, the critical variables for the product and for the manufacturing of the product are determined at **204**. The critical variables are determined by experts in the relevant art, in this case, the semiconductor manufacturing art. As an example, the experts determine that the following factors would be the critical variables for the proposed semiconductor product:

- | | |
|---|--|
| 1. Speed | 2. N2P |
| 3. Vth | 4. Gate Oxide |
| 5. Gate Resistance | 6. R s/d |
| 7. Leff | 8. LDC |
| 9. LDD | 10. S/D Rs (Source/Drain resistance) |
| 11. PolyDelW | 12. Iox |
| 13. RTA sd (rapid thermal anneal s/d) | 14. DopLoss |
| 15. ScreenOx | 16. spacer |
| 17. spacer Thk | 18. spacer Ent |
| 19. PolyThk | 20. G Dopant (gate dopant) |
| 21. s'cide Rs (silicide resistance) | 22. Act Length |
| 23. CD control | 24. SiON Strip (silicon oxynitride strip) |
| 25. G stp | 26. LI bot CD |
| 27. xface s/d Rc | 28. Co QT |
| 29. s'cide pen | 30. G N2.I2 |
| 31. RpdepThk | 32. LI DI CD |
| 33. spacer c/e | 34. s'cide RTA (silicide rapid thermal anneal) |
| 35. Co Thk | 36. XterW (transistor width) |
| 37. VNI/VPI | 38. GateOx QT |
| 39. G R'cle (Gate reticle) | 40. SiON Dep (silicon oxynitride deposition) |
| 41. G Mask QT | 42. Trim Time |
| 43. I2 damage | 44. UDOXthk |
| 45. nitrideDep (nitride deposition) | 46. polishThk (polish thickness) |
| 47. LI etch recipe (local interconnect etch recipe) | 48. LI Thk (local interconnect thickness) |
| 49. LI align (local interconnect alignment) | |
| 50. LI o/e | 51. BMD PCII |
| 52. BMD DepThk | 53. RPE o/e |

Generating such lists of responses and factors is a common part of experimental design practice. At one end of the experimental spectrum are small experiments and short lists of especially important variables (responses and factors). Because there weren't efficient experimental methods to incorporate larger numbers of variables, complex

systems with larger numbers of variables could not be efficiently designed. The present invention provides a method of efficiently designing larger more complex lists of variables. The ability of efficiently designing experimental designs for the larger more complex lists of variables is the major value of the present invention.

5 Using the above-listed critical variables, a classical causal network diagram is created, **Figure 2 Step B, 204**. The terms “cause-effect diagrams,” “causal networks,” and “causal maps” are used in various ways in the relevant art. A cause-effect (or CE) diagram consists of a single response (at the head of a central backbone), major factor groups (as spines branching from the backbone), and lower level factors (off each spine), and perhaps
10 branching off from other low level factors. In modern terms, CE diagrams render causal relationships as a strict hierarchy, with the response as the root node, factor groups as the primary branches and other branches expanding therefrom. Causal networks are generalizations of CE diagrams, with three extensions: (1) the number of responses can be more than one, (2) responses can point causally to other responses, and (3) each factor is
15 represented as one and only one node, even though it may contribute to various other responses and/or to higher-level factors. With newly introduced semiconductor manufacturing processes, transistor characterizations have about $F=50$ factors (F denotes the number of factors of interest) and F_1 is about 8 (F_1 is the number of factors feasible to vary in any single block), the corresponding causal networks are complicated. For this
20 reason, causal networks are transformed into causal maps.

 The transformation of causal networks into causal maps involves the following: (i) the distance between any pair of nodes of a causal network is the minimum number of links of the path connecting them (this is shown below in the internode link-count distance matrix); (ii) the corresponding matrix giving the distances between any pair of nodes is the
25 natural input data structure for multidimensional scaling; and (iii) using a multidimensional scaling algorithm, such as XGvis (Buja et al. 1998) wherein the distance matrix is transformed into node coordinates in $D = 2$ dimensions.

 Causal maps that are constructed in this way have extra information that causal networks do not: (1) a factor closer to a response node plausibly has a stronger effect; (2)
30 two factors close together likely share an interaction; (3) responses sharing many factors cluster; and (4) higher-level factors tend toward the map center.

The causal network diagram created using the critical variables is shown in **Figures 3 & 3A – 3G**. Establishing which critical variables affect other factors creates the causal-effect network. A causal-effect is in the form “cause => effect.” Some of the critical variables are established as “higher-level responses” and are shown enclosed in boxes. Responses shown in **Figure 3** are “Leff” at **302**, “R s/d” at **304** and “Vth(long)” at **306**. The sub-diagrams shown in **Figures 3A – 3G** show some of the critical variables shown in **Figure 3** with additional critical variables “pointing” to them. For example, a critical factor “poly delta W” shown at **308**, **Figure 3** is shown in **Figure 3A** with three arrows from critical variables pointing to it. It should be noted that the critical factor “poly delta W” is therefore a response because other critical variables affect it (the number of arrows that are directed toward it), for example, the critical factor “poly delta W” as shown in **Figure 3A** has 3 arrows directed towards it directly and 5 arrows directed to another critical factor “CD control” **310** and an arrow from the critical factor “CD control” **310** is directed to the critical factor “poly delta W.” Some of the critical variables that are identified as responses, such as the response “poly delta W” are termed “intermediate responses.”

The next step is to create an Internode Link-Count Distance Matrix, **Figure 2, Step C, 208**. **Table 1** is an “internode link-count distance matrix.” The values in the matrix are derived by counting the least number of links from any one node of a factor or response any to any other factor or response in a causal map as shown in **Figure 4**. For example, the internode link-count distance “1” from the factor “speed” to the factor “N2P” is obtained by counting the link from the factor “N2P” to the factor “speed.” Similarly, the internode link-count distance “2” from the factor “speed” to the factor “Vth” is obtained by counting the link from the factor “Vth” to the factor “N2P” and the link from the factor “N2P” to the factor “speed.” The other matrix elements are derived in the same manner.

INTERNODE LINK-COUNT DISTANCE MATRIX
TABLE 1

	1 Speed	2 N2P	3 Vth	4 Gate Ox	5 Gate R	51 BMD PCII	BMD DepThk	RPE o/e
30 Factors (F)								
1 Speed	0	1	2	2	1	4	4	4
2 N2P	1	0	1	1	2	3	3	3
3 Vth	2	1	0	1	3	4	4	4
4 Gate Ox	2	1	1	0	3	4	4	4
35 5 Gate R	1	2	3	3	0	4	4	4
6 R s/d	2	1	2	2	3	2	2	2

	7 Leff	2	1	2	2	3	4	4	4
	8 LDC	3	2	1	2	4	5	5	5
	9 LDD	3	2	3	3	4	3	3	3
	10 S/D Rs	3	2	2	3	2	3	3	3
5	11 PolyDelW	3	2	3	3	4	4	4	4
	12 Iox	4	5	5	6	3	4	4	4
	13 RTAsd	3	2	1	2	3	4	4	4
	14 DopLoss	4	3	3	4	3	4	4	4
	15 ScreenOx	3	4	4	5	2	3	3	3
10	16 spacer	3	2	3	3	4	3	3	3
	17 spacer Thk	4	3	4	4	5	4	4	4
	18 spacer Ent	4	3	4	4	5	4	4	4
	19 PolyThk	4	3	2	3	5	4	4	4
	20 G Dopant	3	2	1	2	4	5	5	5
15	21 s'cide Rs	2	3	3	4	1	3	3	3
	22 Act Length	3	2	3	3	4	3	3	3
	23 CD control								
	24 SiON strip								
	25 G stp								
20	26 LI bot CD								
	27 xface s/d Rc								
	28 Co QT								
	29 s'cide pen								
25	30 G N2.I2								
	31 RpdepThk								
	32 LI DI CD								
	33 spacer c/e								
	34 s'cide RTA								
30	35 Co Thk								
	36 XterW								
	37 VNI/VPI								
	38 GateOx QT								
	39 G R'cle								
35	40 SiON Dep								
	41 G Mask QT								
	42 TrimTime								
	43 I2 damage								
	44 UDOXthk								
40	45 nitrideDep								
	46 polishThk								
	47 LI etch recipe								
	48 LI Thk								
	49 LI align								
	50 LI o/e								
45	51 BMD PCII								
	52 BMD DepThk								
	53 RPE o/e								

etc.

When the Internode Link-Count Distance Matrix is determined, the next step is to apply a multidimensional scaling algorithm to create a D-dimensional (D typically 2 or 3) set of node coordinates called a causal map. Especially when $D = 2$, one can plot the nodes as points on rectangular graph paper, and so both the coordinates themselves and the

resulting graph **400**, **Figure 4** are called causal maps. One example of a multidimensional scaling algorithm is XGvis. Its application to visualizing networks is known and is conventional practice.

The next step is to create a Causal Map, **Figure 2, Step D, 210**. **Figure 4** shows the Causal Map **400** derived from information contained in the causal network shown in **Figures 3 & 3A – 3G** and from information contained in the Internode Link-Count Distance Matrix, **Table 2**.

Once the Causal Map is determined, the next step is to Identify Response Nodes, **Figure 2, Step E, 212**. An inspection of the Causal Map **400**, **Figure 4** shows that the key responses move towards the center and are R s/d, n and p, at **402**, V_{th}, n and p, at **404** and L_{eff}, n and p, at **406**. It is noted that part of the selection of key responses is the exercise of engineering judgment and additional selections or alternate selections could be made.

After the values in **Table 1** are determined and after the Key Responses are determined the next step is to calculate the Map-Based Coefficients a_{ij} , **Figure 2, Step F, 214**. **Table 2** shows the Matrix A (columns V_{th},n to L_{eff},p of **Table 2** make up Matrix A) with values a_{ij} derived from information contained in **Table 1 & Figure 2, Step F, at 214**.

The following example shows how the values in **Table 2** are calculated. Suppose the causal network has factors f1, f2, and f3 all pointing to response r0 having values determined from a similar Internode Link-Count Distance Matrix as shown in **Table 1**.

The portion of the internode link-count distance matrix is:

	r0	f1	f2	f3
r0	0	1	1	1
f1	1	0	2	2
f2	1	2	0	2
f3	1	2	2	0

The causal map coordinates in 2 dimensions (D = 2) are approximately:

	dim1	dim2
r0	0	0
f1	1	0
f2	-0.5	0.866
f3	-0.5	-0.866

Note: r0 is in the middle and f1,f2, and f3 form an equilateral triangle around r0.

The distance between f_i and f_j is square root of 3 (not quite 2), but the distance between r_0 and f_i is 1. The multidimensional scaling would try to balance this, and so the causal map in $D = 2$ might be approximately 1.3 times the above, as shown below:

	dim1	dim2
5		
r_0	0	0
f_1	1.3	0
f_2	-0.65	1.126
f_3	-0.65	-1.126

Note: The distances between f_i and f_j are 2.252 and the distance between r_0 and f_i is 1.3.

The coefficient is determined as follows:

$$(r_0, f_i) = e^{(\alpha * \text{distance}(r_0, f_i)^2)} = e^{(-\alpha * 1.3 * 1.3)} = e^{-1.69}, \text{ since } \alpha \sim 1, \text{ then } (r_0, f_i) = 0.185.$$

The matrix would then be:

	r_0
f_1	0.185
f_2	-0.185 *
f_3	0.185

*The negative sign for factor f_2 is applied by experience and indicates that an increase in the value of factor f_2 produces a decrease in the response r_0 .

TABLE 2
Map Based Coefficients (Matrix A)

	knob	Vth,n	R s/d,n	Leff,n	Vth,p	R s/d,p	Leff,p
	1 XterW		-0.107888			-0.107888	
	2 I2 damage		0.025962				
	3 VNI	-0.221786					
30	4 VPI				0.221786		
	5 GateOx QT	0.137194		-0.137194			
	6 Gate Ox	0.246187		-0.246187			
	7 PolyThk	0.116432	0.119731		-0.116432	0.119731	
	8 G N2.I2	-0.052234		-0.104468			
35	9 SiOn Dep			0.065866			0.065866
	10 G Mask QT			0.060969			0.060969
	11 G stp			0.158818			0.158818
	12 SiON strip			-0.064624			-0.064624
	13 LDC n	-0.179057		-0.196067			
40	14 LDC p				0.179057		-0.196067
	15 LDD n		-0.227165	-0.166824			
	16 LDD p					-0.227165	-0.166824
	17 UDOX1		0.082092			0.082092	
	18 Iox		0.025134			0.025134	
45	19 spacer Thk		0.108862	0.109174		0.108862	0.109174
	20 spacer o/e		-0.152447	-0.108867		-0.152447	-0.108867
	21 spacEtchEnt						
	22 UDOX2		0.026926			0.026926	
	23 RpdepThk		-0.10924			-0.10924	
50	24 RTAsd	-0.202057	-0.113617	-0.069277	0.202057	-0.113617	-0.069277
	25 RPE o/e		-0.107437			-0.107437	

	26 Co QT	0.109251	0.109251
	27 Co Thk	-0.113116	-0.113116
	28 s'cide RTA1	-0.116867	-0.116867
	29 s'cide RTA2	-0.116867	-0.116867
5	30 LI Thk	-0.096417	-0.096417
	31 polish Thk	0.09219	0.09219
	32 LI DICD	-0.87249	-0.087249
	33 LI alignX	0.078742	0.078742
	34 LI alignY	0.078742	0.078742
10	35 LI o/e	0.106432	0.106432
	36 BMDDepTh	0.101266	0.101266
	37 BMD PCII	0.101964	0.101964

Note: Blanks (or zeros) indicate that a particular critical factor does not affect the particular response, i.e., there is no "cause-effect."

After the map-based coefficients (Matrix **A**) are determined, **Figure 2, Step F, 214**, k is set to 0 and the matrix U_k =(null, zero-row matrix), **Figure 2, Step G, 216**.

An initial base design matrix, **X**, is established by statistician experts in the field, **Figure 2, Step H, 218**. The base design matrix **X** shows in a generic way the form of the design; it is generic in the sense that particular factors are not yet assigned to particular **X**-columns.

Table 3 show two design alternatives, one design alternative is a "split-plot only design" and the other is an "interblock" design. The split-plot design varies factors only within the lot (block). The interblock design varies factors both from lot to lot and within lot. The former set of factors are conventionally termed skew factors, the latter are termed split factors. In each of the two design alternatives L_8 and L'_8 are matrices as follows:

$$L_8 = \begin{bmatrix} - & - & - & - \\ + & - & - & + \\ - & + & - & + \\ + & + & - & - \\ - & - & + & + \\ + & - & + & + \\ - & + & + & - \\ + & + & + & + \end{bmatrix} \quad \text{and } L'_8 = \begin{bmatrix} - & - & - & + \\ + & - & - & - \\ - & + & - & - \\ + & + & - & - \\ - & - & + & - \\ + & - & + & + \\ - & + & + & + \\ + & + & + & - \end{bmatrix}$$

5

TABLE 3

	split-plot design	interblock design
10	$\begin{bmatrix} 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ [L_8] & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \end{bmatrix}$	$\begin{bmatrix} +0-0 & +0-- & +-0+ \\ +0-0 & +0-- & +-0+ \\ +0-0 & +0-- & +-0+ \\ [L_8] & +0-- & +-0+ \\ +0-0 & +0-- & +-0+ \\ +0-0 & +0-- & +-0+ \\ +0-0 & +0-- & +-0+ \\ +0-0 & +0-- & +-0+ \end{bmatrix}$
15		
20	$\begin{bmatrix} 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & [L_8] & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \end{bmatrix}$	$\begin{bmatrix} +-0+ & -+00 & 0+- \\ +-0+ & -+00 & 0+- \\ +-0+ & -+00 & 0+- \\ [L_8] & -+00 & 0+- \\ +-0+ & -+00 & 0+- \\ +-0+ & -+00 & 0+- \\ +-0+ & -+00 & 0+- \\ +-0+ & -+00 & 0+- \end{bmatrix}$
25		
30	$\begin{bmatrix} 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & [L'_8] \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \end{bmatrix}$	$\begin{bmatrix} 0+- & -+0+ & -+0+ \\ 0+- & -+0+ & -+0+ \\ 0+- & -+0+ & -+0+ \\ 0+- & -+0+ & [L'_8] \\ 0+- & -+0+ & -+0+ \\ 0+- & -+0+ & -+0+ \\ 0+- & -+0+ & -+0+ \\ 0+- & -+0+ & -+0+ \end{bmatrix}$
35		
40	$\begin{bmatrix} 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & 0000 \\ 0000 & 0000 & [L'_8] \end{bmatrix}$	$\begin{bmatrix} -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \\ -0+0 & 0-+- & 0-++ \end{bmatrix}$

45 P is a permutation matrix and the matrix product \mathbf{XP} constitutes a rearrangement of the columns \mathbf{X} . The matrix \mathbf{A} is a matrix of coefficients linking (linearly) the full 37 dimensional space to the 6 dimensional intermediate variable space, (37 factors (first column) and 6 responses (column headings) **Table 2**). The established initial design matrix \mathbf{X} does not associate particular factors with particular columns of \mathbf{X} , but it does

50 describe the overall patterns of the experiment, that is, how many split factors per block, how many skew factors per block, how many runs (wafers) per block. A block is a set of experimental runs (wafers) processed together (in semiconductor processing; a lot). Split factors are factors that take on at least two values within a given lot (block). Skew factors

are factors that take on different values from nominal (0), but are constant within the block.

For any particular assignment of X-columns to factors one can compute a score. This score is related to a prediction of how spread out the eventual results (responses) measured on the wafers (runs, rows of X) are predicted to be. The better designs will have better (higher) scores. This number is calculated by assigning particular columns of X to particular factors, one-to-one, with no duplicate assignments and no factors left out.

A list-to-list assignment, with a score about which assignments are worse or better, is known as “the traveling salesman problem.” In the conventional description of the traveling salesman problem, a salesman needs to visit a list of cities once each and wants to minimize the driving time. In that context, a salesman has one list—of cities—and another list, the number 1, 2, 3, 4, and needs to assign each number to each city. Assigning 1 to city B means visit that city first, 2 to city D means visit that city second, and so on. Such an algorithm gives an X-column-to-factor assignment with the biggest score of any assignment considered. The R step/E step algorithm discussed below is one particular traveling-salesman-solving algorithm.

Conceptually, if an assignment is made of X-columns and factors at random and if another random assignment is made, there is a 50% chance that a higher score would be obtained. If a third random assignment is made, there is a chance of 33% of achieving a higher score than the previous two assignments, etc. As can be appreciated, random guessing (R step) bogs down and offers no improvement after numerous tries. What is done is pick the best so far and see if, swapping just one pair of factors' X-columns achieves a higher score. Doing this one pair at a time for all pairs of factors is the done by the E algorithm discussed below and is repeated until all pairs have been examined without an improvement.

The optimal design algorithm including the algorithm R(k) and the algorithm E(k) are run with $k = k + 1$. **Figure 2, Step I, 220.** $Y(P) = (I - B(B^T B)^{-1} B^T)[(X P) / U]A$ & **Wynn's criterion** is run. Note that the term $(I - B(B^T B)^{-1} B^T)$ is a projection matrix that acts to filter out effects correlated with **B**, leaving only effects not correlated with **B**, that is, it serves as a blocking function. **B** is a matrix of dummy variables representing the blocks. After $Y(P)$ is defined, the algorithms R(k) and E(k) are run. R(k) is an algorithm

to choose the best of n_r random column permutation matrices \mathbf{P} . $E(k)$ is an algorithm is to choose the best column permutation matrix \mathbf{P} that is “near” a previous solution. (On the first iteration of algorithm E, the previous solution comes from algorithm R; thereafter, the previous solution comes from the last application of algorithm E.) Note that \mathbf{A}/\mathbf{B} means:

5 “append the rows of \mathbf{B} to the rows of \mathbf{A} , with the rows of \mathbf{A} on top.

Wynn’s criterion: If $\{y_i: i = 1 \dots n\}$ denotes a set of k -dimensional points comprising a possible experimental design, and the $n \times n$ matrix \mathbf{C} is defined with typical element $c(i,j) = \exp(-\|y_i - y_j\|^2)$. Wynn’s criterion is that a design $\{y_i\}$ is better when $\det \mathbf{C}$ is larger (“det” denotes the matrix determinant). If \mathbf{C} is interpreted as a correlation matrix, 10 $\det(\mathbf{C})$ represents the generalized variance, which Wynn’s criterion maximizes. This is achieved by moving the points $\{y_i: i=1 \dots n\}$ far from one another.

\mathbf{C} , such that $C(i,j) = \exp\{-\|y_i - y_j\|^2\}$. Distances among points $\{y_i\}$ are transformed into “correlations” as discussed above. The points $\{y_i\}$ are the rows of matrix $\mathbf{Y}(\mathbf{P})$.

Algorithm R (k):

15 Let \mathbf{X}' denote the best current design, with criterion value c' , which is a scalar and initially $c' = -\infty$.

Loop in $r=1$ to n_R

Randomly rearrange columns of base design matrix \mathbf{X} , resulting in \mathbf{X}_r

Form \mathbf{W}_r by adding to matrix \mathbf{X}_r the rows of \mathbf{U}_{k-1}

20 $\mathbf{Y}_{r1} = \mathbf{W}_r \mathbf{A}$

$\mathbf{Y}_{r2} = [\mathbf{I} - \mathbf{B}(\mathbf{B}^T \mathbf{B})^{-1} \mathbf{B}^T] \mathbf{Y}_{r1}$

$c_r = \det \mathbf{C}(\mathbf{Y}_{r1}) \times \det \mathbf{C}(\mathbf{Y}_{r2})$

if it is determined at 222, Figure 2, that $c_r > c'$ then $\mathbf{X}' = \mathbf{X}_r$ and $c' = c_r$

end Loop r .

25 The function of algorithm R is to choose the best of n_r random column permutation matrices \mathbf{P} .

Algorithm E(k):

Let \mathbf{X}' denote the best current design, with criterion value c' , the scalar from algorithm R.

30 Label E: NoImprovement=true

$\mathbf{X}'' = \mathbf{X}'$

$c'' = c'$

Loop in $i=1$ to $n-1$

Loop in $j = i+1$ to n
 Form matrix X_{ij} by swapping columns i, j of X'
 Form W_{ij} by adding to matrix X_{ij} the rows of U_{k-1} *
 $Y_{e1} = W_{ij}A$
 5 $Y_{e2} = [I - B(B^T B)^{-1} B^T] Y_{e1}$
 $c_{ij} = \det C(Y_{e1}) \times \det C(Y_{e2})$
 If $c_{ij} > c''$, then $X'' = X_{ij}$ and $c'' = c_{ij}$, NoImprovement = false
 end Loop j
 end Loop i
 10 if NoImprovement then STOP else go to Label E.

Note 1: If a factor with a strong factor is split, Y_{r1} & Y_{r2} are both more dispersed, hence larger criterion. If some factor is skewed instead, Y_{r1} is dispersed, but not Y_{r2} . In this way this criterion gives double points for splitting strong factors.

* Note 2: It is conventional practice for statisticians to term vertical elements as
 15 "columns" and the horizontal elements as "rows."

Tables 4 – 6 are examples of designs and solutions for particular 18-wafer lots
 and were extracted from a larger solution of 6-8 lots. Each design and solution, for
 example, the design and solution in Table 4 is a design U_k indicated at 224, Figure 2,
 20 Step J.

After the optimal design algorithm is run Step I, Figure 2, 220 is run it is
 determined at 222 if the design U_k is large enough. If it is determined at 220 that there the
 design U_k is not large enough the flow returns to Step I and the optimal design algorithm is
 repeated until it is determined at 222 that the design U_k is large enough. When this occurs
 25 the flow goes to Step J and the design experiment U_k is complete and the flow returns to
 Figure 1, Step D as indicated by arrow 226.

When the flow returns to Step D, Figure 1, prototype wafers using the designed
 experiment are manufactured. As discussed above, Tables 4 – 6 each show 1 lot (18
 wafers) of a solution and design that could be 6-8 wafers. The manufactured prototype
 30 wafers are tested and model responses are measured at 114, Figure 1, Step E.

It is determined at 116 whether the Model Responses are adequate. If the model
 responses are not adequate, the flow returns to Step I, Figure 2 as indicated by arrow 118.
 Step I, Figure 2 is the step of running the Optimal design algorithm and Steps I, Figure 2

through **Step E, Figure 1** are repeated until it is determined at **116, Figure 1** that the model is adequate.

After the model is determined to be adequate at **116**, the tolerances are assessed/proposed at **Step F, Figure 1, 120**. Whether the proposed tolerances are manufacturable is determined at **122, Figure 1** and if the proposed tolerances are not manufacturable, the flow returns to **Step B** as indicated by arrow **126**. The development of the idea then begins anew at **Step B**. If it is determined at **122, Figure 1** that the proposed tolerances are manufacturable, the best solution of the design is sent to production at **Step G, Figure 1, 124**.

Table 4 is a lot (block) of 18 wafers and the total solution could be 6 – 8 lots. After a total number of lots is completed, the wafers are manufactured and tested to validate the experiment and to choose the best set of factors to use for full scale production.

Table 4 is an example Split Sheet for one of 6 lots and is a part of a matrix **[XP]** and shows the results of running the optimization algorithm. A split sheet includes instructions for processing the wafers of one particular lot (one block). Format is **(XP)^T**: that is, the columns are wafers in a split sheet, and the rows are factors. (Conventional experimental design notation has a reverse convention: the rows (the vertical elements) are wafers and the columns (the horizontal elements) are factors.)

20

TABLE 4

wafer # →		1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18
Factor																			
	1 XterW	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
25	2 I2 damage	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0 <=**
	=>*3 VNI	0	-	+	+	-	-	+	+	-	+	-	-	+	+	-	-	+	0
	4 VPI	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	5 GateOx QT	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	6 GateOx	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
30	7 PolyThk	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	=> 8 G N2 I2	0	-	-	+	+	-	-	+	+	-	-	+	+	-	-	+	+	0
	9 SiOn Dep	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	10 G Mask QT	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	11 G stp	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
35	12 SiOn strip	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	=> 13 LDC n	0	-	-	-	-	-	-	-	-	+	+	+	+	+	+	+	+	0
	14 LDC p	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	15 Iox	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	16 LDD n	0	-	+	-	+	-	+	-	+	-	+	-	+	-	+	-	+	0
40	17 LDD p	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=

	18 UDOX1	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	19 spacer Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	20 spacer o/e	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0
5	=> 21 spacerEtchEnt	0	-	+	+	-	+	-	-	+	-	+	+	-	+	-	+
	22 UDOX2	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	23 RpdepThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	=> 24 RTAsd	0	-	-	-	-	++	+	+	-	-	-	+	+	+	+	+
	25 RPE o/e	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	26 Co QT	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
10	27 Co Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	28 s'cide RTA1	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	29 s'cide RTA2	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	30 LI Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	31 polishThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
15	32 LI DICD	0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0<=
	33 LI alignX	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	34 LI alignY	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	35 LI o/e	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	36 BMD DepThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
20	37 BMD PCII	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0

* The left-side arrows indicate, "split factors."

** The right-side arrows mean that the factors varied in this lot for lot-to-lot experiment (skew factors).

Note: In Tables 4 – 6, the "0" means that the factor has not been increased or decreased from the nominal value, that the "-" means that the factor has been decreased and that the "+" means that the factor has been increased.

Table 5 is a Split Lot with No Skewed Factors

		TABLE 5																	
wafer # →		1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18
35	Factor																		
	1 XterW	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	2 I2 damage	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	3 VNI	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
40	4 VPI	0	-1	-1	-1	-1	-1	-1	-1	-1	1	1	1	1	1	1	1	1	0
	5 GateOx	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	6 GateOx	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	7 PolyThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
45	8 G N2 I2	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	9 SiOn Dep	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	10 G Mask QT	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	11 G stp	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
50	12 SiOn strip	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	13 LDC n	0	-1	-1	-1	-1	1	1	1	1	-1	-1	-1	-1	-1	1	1	1	0
	14 LDC p	0	-1	1	-1	1	1	-1	1	-1	1	-1	1	-1	-1	1	-1	1	0
	15 Iox	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	16 LDD n	0	-1	1	-1	1	-1	1	-1	1	-1	1	-1	1	-1	1	-1	1	0
	17 LDD p	0	-1	-1	1	1	1	-1	-1	1	1	-1	-1	-1	-1	-1	1	1	0
	18 UDOX1	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	19 spacer Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0

	20 spacer o/e	0	-1	-1	1	1	-1	-1	1	1	-1	-1	1	1	0
	21 spacerEtchEnt	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	22 UDOX2	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	23 RpdepThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0
5	24 RTAsd	0	-1	1	1	-1	1	1	-1	1	1	-1	-1	1	0
	25 RPE o/e	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	26 Co QT	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	27 Co Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	28 s'cide RTA1	0	0	0	0	0	0	0	0	0	0	0	0	0	0
10	29 s'cide RTA2	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	30 LI Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	31 polishThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	32 LI DICD	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	33 LI alignX	0	0	0	0	0	0	0	0	0	0	0	0	0	0
15	34 LI alignY	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	35 LI o/e	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	36 BMD DepThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	37 BMD PCII	0	0	0	0	0	0	0	0	0	0	0	0	0	0

20

Table 6 is a Split Lot that includes skewed factors.

		TABLE 6																	
wafer # →		1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18
25	1 XterW	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	2 I2 damage	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	3 VNI	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	4 VPI	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	5 GateOx QT	0	-1	-1	1	1	-1	-1	1	1	-1	-1	1	1	-1	-1	1	1	0
30	6 GateOx	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	7 PolyThk	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	8 G N2 I2	0	-1	-1	-1	-1	1	1	1	1	-1	-1	-1	-1	1	1	1	1	0
	9 SiOn Dep	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
35	10 G Mask QT	0	-1	1	-1	1	-1	1	-1	1	-1	1	-1	1	-1	1	-1	1	0
	11 G stp	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	12 SiOn strip	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	13 LDC n	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	14 LDC p	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
40	15 Iox	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	16 LDD n	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	17 LDD p	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	18 UDOX1	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	19 spacer Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
45	20 spacer o/e	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	21 spacerEtchEnt	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	22 UDOX2	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	23 RpdepThk	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	24 RTAsd	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
50	25 RPE o/e	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	26 Co QT	0	-1	-1	-1	-1	-1	-1	-1	-1	1	1	1	1	1	1	1	1	0
	27 Co Thk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	28 s'cide RTA1	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	29 s'cide RTA2	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
55	30 Li Thk	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0
	31 polishThk	0	-1	1	1	-1	1	-1	-1	1	1	-1	1	-1	-1	-1	-1	1	0

32 LI DICD	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
33 LI alignX	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
34 LI alignY	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
35 LI o/e	0	-1	1	1	-1	-1	1	1	-1	1	-1	1	1	-1	-1	1
5 36 BMD DepThk	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
37 BMD PCII	0	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	-1	0

In summary, the described invention thus provides a method for a systematic approach to forming experimental designs for large, complex systems after an idea for a product is formed.

The foregoing description of the embodiment of the invention has been presented for purposes of illustration and description. It is not intended to be exhaustive or to limit the invention to the precise form disclosed. Obvious modifications or variations are possible in light of the above teachings. The embodiment was chosen and described to provide the best illustration of the principles of the invention and its practical application to thereby enable one of ordinary skill in the art to utilize the invention in various embodiments and with various modifications as are suited to the particular use contemplated. All such modifications and variations are within the scope of the invention as determined by the appended claims when interpreted in accordance with the breadth to which they are fairly, legally, and equitably entitled.